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LOGINID: SSSPTA1623PAZ

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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Welcome to STN International
NEWS
                 Web Page URLs for STN Seminar Schedule - N. America
                 "Ask CAS" for self-help around the clock
NEWS
NEWS
     3 May 12
                 EXTEND option available in structure searching
        May 12
NEWS
                 Polymer links for the POLYLINK command completed in REGISTRY
NEWS
     5 May 27
                 New UPM (Update Code Maximum) field for more efficient patent
                 SDIs in CAplus
NEWS
     6 May 27
                 CAplus super roles and document types searchable in REGISTRY
NEWS
     7
         Jun 28
                 Additional enzyme-catalyzed reactions added to CASREACT
NEWS
         Jun 28
                 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG,
                 and WATER from CSA now available on STN(R)
NEWS
        Jul 12
                 BEILSTEIN enhanced with new display and select options,
                 resulting in a closer connection to BABS
NEWS 10
        Jul 30
                 BEILSTEIN on STN workshop to be held August 24 in conjunction
                 with the 228th ACS National Meeting
NEWS 11
                 IFIPAT/IFIUDB/IFICDB reloaded with new search and display
         AUG 02
                 fields
NEWS 12
        AUG 02
                 CAplus and CA patent records enhanced with European and Japan
                 Patent Office Classifications
        AUG 02
NEWS 13
                 STN User Update to be held August 22 in conjunction with the
                 228th ACS National Meeting
NEWS 14
         AUG 02
                 The Analysis Edition of STN Express with Discover!
                 (Version 7.01 for Windows) now available
NEWS 15
        AUG 04
                 Pricing for the Save Answers for SciFinder Wizard within
                 STN Express with Discover! will change September 1, 2004
NEWS EXPRESS
              JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT
              MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
              AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
              STN Operating Hours Plus Help Desk Availability
NEWS HOURS
NEWS INTER
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              Welcome Banner and News Items
              Direct Dial and Telecommunication Network Access to STN
NEWS PHONE
NEWS WWW
              CAS World Wide Web Site (general information)
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FILE 'HOME' ENTERED AT 11:57:09 ON 18 AUG 2004

=> file reg\

'REG\' IS NOT A VALID FILE NAME

SESSION CONTINUES IN FILE 'HOME'

Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> file reg

COST IN U.S. DOLLARS SINCE FILE TOTAL

FULL ESTIMATED COST ENTRY SESSION 0.21 0.21

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STRUCTURE FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2 DICTIONARY FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>	e	trip.	hosgene/	cn

E1	1 TRIPHOS (FEED SUPPLEMENT)/CN
E2	1 TRIPHOS-I/CN
E3	1> TRIPHOSGENE/CN
E4	1 TRIPHOSGENE-PHENYLPHOSPHONIC DICHLORIDE-1,1-BIS(4-HYDROXYPHE
	NYL)CYCLOHEXANE COPOLYMER/CN
E5	1 TRIPHOSPHABORETANE/CN
E6	1 TRIPHOSPHADEN/CN
E7	1 TRIPHOSPHAGERMETANE/CN
E8	1 TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIETH
	YL-/CN
E9	1 TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIPHE
	NYL-/CN
E10	1 TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIPHE
	NYL-, (1A,2B,3A)-/CN
E11	1 TRIPHOSPHAGERMETANE, 4,4-DICHLORO-1,2,3-TRIS(1,1-DIMETHYLETH
	YL)-/CN
E12	1 TRIPHOSPHAMIDE/CN

=> e3

L1 1 TRIPHOSGENE/CN

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 4.85 5.06

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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 11

854 L1 L2

=> chloroformate

18822 CHLOROFORMATE 1670 CHLOROFORMATES

19404 CHLOROFORMATE L3

(CHLOROFORMATE OR CHLOROFORMATES)

=> 12 and 13

141 L2 AND L3

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION 8.35

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 12:00:06 ON 18 AUG 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> file reg
COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 0.46 8.81

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STRUCTURE FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2 DICTIONARY FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

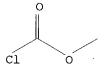
Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>

Uploading C:\Examination Auxillary files\10665410\10665410 chloroformate core.str



5

chain nodes:
1 2 3 4 5
chain bonds:
1-2 2-3 2-5 3-4
exact/norm bonds:
2-3 2-5 3-4
exact bonds:
1-2

Match level:

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS

L5 STRUCTURE UPLOADED

=> d 15 L5 HAS NO ANSWERS

L5 HAS NO ANSWERS
L5 STR

c1

Structure attributes must be viewed using STN Express query preparation.

=> search 15 sss ful FULL SEARCH INITIATED 12:00:39 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 5459 TO ITERATE

100.0% PROCESSED 5459 ITERATIONS SEARCH TIME: 00.00.01

2680 ANSWERS

L6 2680 SEA SSS FUL L5

=> file caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 155.42 164.23

FULL ESTIMATED COST

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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 16/prep

18704 L6 3184864 PREP/RL

L7 1571 L6/PREP

(L6 (L) PREP/RL)

=> d his

(FILE 'HOME' ENTERED AT 11:57:09 ON 18 AUG 2004)

FILE 'REGISTRY' ENTERED AT 11:57:45 ON 18 AUG 2004 E TRIPHOSGENE/CN

L1 1 E3

FILE 'CAPLUS' ENTERED AT 11:58:08 ON 18 AUG 2004

L2 854 L1

L3 19404 CHLOROFORMATE

L4 141 L2 AND L3

FILE 'CAPLUS' ENTERED AT 12:00:06 ON 18 AUG 2004

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L6
             2680 SEARCH L5 SSS FUL
      FILE 'CAPLUS' ENTERED AT 12:00:46 ON 18 AUG 2004
L7
             1571 L6/PREP
=> 12 and 17
              38 L2 AND L7
=> amine
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          235258 AMINES
L9
          379582 AMINE
                    (AMINE OR AMINES)
=> 19 and 18
               5 L9 AND L8
=> d 110 1-5 ti fbib abs
L10 ANSWER 1 OF 5 CAPLUS COPYRIGHT 2004 ACS on STN
      Preparation of 5-membered heterocycle derivatives for treating
      neurodegenerative disorders or pain
      2004:550745 CAPLUS
ΑN
DN
      141:106475
      Preparation of 5-membered heterocycle derivatives for treating
      neurodegenerative disorders or pain
      Chabrier De Lassauniere, Pierre-Etienne; Harnett, Jeremiah; Bigg, Dennis;
IN
      Liberatore, Anne-Marie; Pommier, Jacques; Lannoy, Jacques; Thurieau,
      Christophe
PA
SO
     U.S. Pat. Appl. Publ., 150 pp., Cont.-in-part of U.S. Ser. No. 89,993.
      CODEN: USXXCO
DΤ
     Patent
     English
LA
FAN.CNT 2
      PATENT NO.
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                                      DATE
                                                    APPLICATION NO.
                                                                                DATE
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FR 2000-10151

A 20000801

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FR 2000-11169 A 20000901
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PATENT FAMILY INFORMATION:
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                            KIND
                                      DATE APPLICATION NO.
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                                                                                DATE
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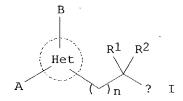
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                                      20040708
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FR	2000-10151	A	20000801
FR	2000-11169	А	20000901
WO	2000-FR2805	W	20001010
JΡ	1989-4943	Α	20010410
JP	1990-1811	A	20020214
US	2002-89993	` A2	20020404

GΙ



AΒ The invention relates to thiazole, oxazole, imidazole, isoxazole and isoxazoline derivs. of general formula (I) [wherein Het = thiazole, oxazole, imidazole, isoxazole or isoxazoline; n = an integer from 0 to 6; A = optionally substituted aromatic radical; B = H, alkyl, Ph; R1, R2 = H, alkyl, cycloalkyl; Ω = NR46R47 or OR48; R46, R47 = H, alkyl, cycloalkyl, (CH2)k-CO2R51; R51 = alkyl, haloalkyl; R48 = H, alkyl]. compds. have advantageous pharmacol. properties which allow their use in a medicament intended to inhibit monoamine oxidases (MAO) and/or lipidic peroxidn. and/or to act as modulators of the sodium channels and notably their use in therapeutics for treating (1) central or peripheral nervous system, (2) neurodegenerative disorders selected from Parkinson's disease, Alzheimer's disease, Huntington's chorea and amyotrophic lateral sclerosis or (3) pain selected from the group consisting of postoperative pain. migraine, neuropathic pain, central pain, chronic inflammatory pain and pain linked to a cancer. Thus, 2-[[[(1,1-dimethylethoxy)carbonyl]methyl]a mino]ethanethioamide (4.3 q, 2.11 mmol) and 2-bromo-1-(3,5-di-tert-butyl-4hydroxyphenyl)ethanone (6,9 g, 2,11 mmol) were dissolved in 75 mL benzene under argon atmospheric and stirred at ambient temperature for 12 h to give,

after

workup and silica gel chromatog., $4-[3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl]-N-[(1,1-dimethylethoxy)carbonyl]-N-methyl-2-thiazolemethanamine which was treated with CF3CO2H and triethylsilane in 50 mL CH2Cl2 to give, after workup, <math>4-[3,5-bis(1,1-dimethylethyl)-4-hydroxyphenyl]-N-methyl-2-thiazolemethanamine (II). II showed IC50 of lower than 10 <math display="inline">\mu M$ for inhibiting lipid peroxidn. of the cerebral cortex of rats.

- L10 ANSWER 2 OF 5 CAPLUS COPYRIGHT 2004 ACS on STN
- TI Improved synthesis of (dialkylamino)pyrrolines
- AN 2003:550006 CAPLUS
- DN 139:364788
- TI Improved synthesis of (dialkylamino)pyrrolines
- AU Flosser, David A.; Olofson, Roy A.
- CS Department of Chemistry, The Pennsylvania State University, University Park, PA, USA
- SO Synthetic Communications (2003), 33(12), 2045-2052 CODEN: SYNCAV; ISSN: 0039-7911
- PB Marcel Dekker, Inc.
- DT Journal
- LA English
- OS CASREACT 139:364788

GΙ

AB The title compds. (I; R = NBu2, piperidino, NEt2) were prepared in 80-94% yield by reaction of I (R = OMe) with amines and their hydrochlorides. In initial assays, the pyrrolinium salts obtained on alkylation of I (R = NBu2) are excellent "naked halide" catalysts.

RE.CNT 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 3 OF 5 CAPLUS COPYRIGHT 2004 ACS on STN

TI Novel chiral surfactant for the separation of enantiomers by micellar electrokinetic capillary chromatography

AN 1995:116456 CAPLUS

DN 122:71192

TI Novel chiral surfactant for the separation of enantiomers by micellar electrokinetic capillary chromatography

AU Mazzeo, Jeffrey R.; Grover, Edward R.; Swartz, Michael E.; Petersen, John S.

CS Waters Chromatography, 34 Maple Street, Milford, MA, 01757, USA

SO Journal of Chromatography, A (1994), 680(1), 125-35 CODEN: JCRAEY; ISSN: 0021-9673

DT Journal

LA English

An novel chiral surfactant was prepared as both enantiomeric forms, (R)— and (S)—N—dodecoxycarbonylvaline, and employed for the separation of enantiomeric mixts. by micellar electrokinetic capillary chromatog. (MECC). The enantioselectivities (α) obtained for twelve typical pharmaceutical amines using the (S)—surfactant were compared to those obtained with (S)—N—dodecanoylvaline, a chiral surfactant described in the literature. Higher enantioselectivities were seen for ten of the twelve compds. using (S)—N—dodecoxycarbonylvaline. Furthermore, (S)—N—dodecoxycarbonylvaline had significantly less background absorbance in the low UV. It is shown that exact enantiomer migration order reversal can be obtained by individually employing both enantiomeric forms of the surfactant. For ionizable compds. like the amines examined here, enantioselectivity can be optimized by changing the pH of the MECC buffer. Partitioning is optimized through surfactant concentration, organic additives

and

pH. The ability to achieve fast chiral sepns. is shown. A separation of ephedrine enantiomers in urine is shown, with the only sample preparation being filtration.

L10 ANSWER 4 OF 5 CAPLUS COPYRIGHT 2004 ACS on STN

TI Triphosgene: a crystalline phosgene equivalent

AN 1987:575482 CAPLUS

DN 107:175482

TI Triphosgene: a crystalline phosgene equivalent

AU Eckert, Heiner; Forster, Barbara

CS Org. Chem. Inst. Tech. Univ. Muenchen, Garching, D-8046, Fed. Rep. Ger.

SO Angewandte Chemie (1987), 99(9), 922-3 CODEN: ANCEAD; ISSN: 0044-8249

DT Journal

LA German

OS CASREACT 107:175482

AB (C13CO)2CO (I) was used for chloroformylation, carbonylation, chlorination, and dehydration. Thus, when treated with I, C13CCMe2OH gave 91% C13CCMe2O2CCl, o-MeC6H4NH2 gave 82% o-MeC6H4NCO, PhCH2CO2H gave 11%

PhCH2COC1, and RCH2CH2NHCHO (R = morpholino) gave 74% RCH2CH2NC.

L10 ANSWER 5 OF 5 CAPLUS COPYRIGHT 2004 ACS on STN

TI Bis(trichloromethyl) carbonate as an alternative reagent for phosgene

AN 1987:4294 CAPLUS

DN 106:4294

TI Bis(trichloromethyl) carbonate as an alternative reagent for phosgene

IN Eckert, Heiner

PA Fed. Rep. Ger.

SO Ger. Offen., 17 pp.

CODEN: GWXXBX

DT Patent

LA German

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ΡI	DE 3440141	A1	19860507	DE 1984-3440141 DE 1984-3440141	19841102 19841102
				DE 1904-3440141	19041102

OS CASREACT 106:4294

GΙ

1'_

$$H_2N$$
 CH_2 NH_2 II OCN CH_2 NCO

AB (Cl3CO)2CO (I), prepared by chlorination of (MeO)2CO, is used as an alternative to COCl2 in, e.g., the preparation of isocyanates, diisocyanates, chloroformates, and polycarbonates, etc., which find use as intermediates for plastics, pharmaceuticals, herbicides, and insecticides, etc. Thus, 5.95 g diamine II and 5.94 g I in o-Cl2C6H4 were heated at 170° for 3 h to give 84% diisocyanate III.

=> 99999file reg

0 99999FILE

1060 REG

50 REGS

1105 REG

(REG OR REGS)

L11

0 99999FILE REG

(99999FILE(W)REG)

=> d 11

YOU HAVE REQUESTED DATA FROM FILE 'REGISTRY' - CONTINUE? (Y)/N:n

=> file reg

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	33.61	197.84
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-3.68	-3.68

FILE 'REGISTRY' ENTERED AT 12:08:19 ON 18 AUG 2004

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STRUCTURE FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2 DICTIONARY FILE UPDATES: 17 AUG 2004 HIGHEST RN 727974-89-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> d 11

CI

COM

L1ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN RN 32315-10-9 REGISTRY CN Methanol, trichloro-, carbonate (2:1) (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES: Carbonic acid, bis(trichloromethyl) ester (6CI, 8CI) OTHER NAMES: CN Bis(trichloromethyl) carbonate CN Triphosgene FS 3D CONCORD MF C3 C16 O3

LCSTN Files: BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS, CASREACT,

CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, GMELIN*, HODOC*, MEDLINE, MSDS-OHS, PROMT, SYNTHLINE, TOXCENTER, USPAT2, USPATFULL

(*File contains numerically searchable property data) Other Sources: EINECS**, TSCA**

(**Enter CHEMLIST File for up-to-date regulatory information)

CAplus document type: Conference; Journal; Patent

RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); PROC (Process); RACT (Reactant or reagent); USES (Uses)

Roles for non-specific derivatives from patents: PREP (Preparation); PROC (Process); RACT (Reactant or reagent); USES (Uses)

Roles from non-patents: ANST (Analytical study); FORM (Formation, nonpreparative); MSC (Miscellaneous); OCCU (Occurrence); PREP (Preparation); PROC (Process); PRP (Properties); RACT (Reactant or reagent); USES (Uses)

RLD.NP Roles for non-specific derivatives from non-patents: PREP (Preparation); PRP (Properties); RACT (Reactant or reagent)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

848 REFERENCES IN FILE CA (1907 TO DATE)

10 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

854 REFERENCES IN FILE CAPLUS (1907 TO DATE)

3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> logoff hold COST IN U.S. DOLLARS

SINCE FILE TOTAL

SINCE FILE

FULL ESTIMATED COST

ENTRY SESSION 3.03 200.87

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

TOTAL

ENTRY SESSION

CA SUBSCRIBER PRICE

0.00 -3.68

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 12:10:13 ON 18 AUG 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * * SESSION RESUMED IN FILE 'REGISTRY' AT 12:32:04 ON 18 AUG 2004 FILE 'REGISTRY' ENTERED AT 12:32:04 ON 18 AUG 2004 COPYRIGHT (C) 2004 American Chemical Society (ACS)

COST IN U.S. DOLLARS

SINCE FILE

TOTAL SESSION

FULL ESTIMATED COST

ENTRY 3.03

200.87

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

CA SUBSCRIBER PRICE

ENTRY 0.00

SESSION -3.68

=> d his

(FILE 'HOME' ENTERED AT 11:57:09 ON 18 AUG 2004)

FILE 'REGISTRY' ENTERED AT 11:57:45 ON 18 AUG 2004 E TRIPHOSGENE/CN

L1 1 E3

FILE 'CAPLUS' ENTERED AT 11:58:08 ON 18 AUG 2004

L2 854 L1

L3 19404 CHLOROFORMATE

L4 141 L2 AND L3

FILE 'CAPLUS' ENTERED AT 12:00:06 ON 18 AUG 2004

FILE 'REGISTRY' ENTERED AT 12:00:13 ON 18 AUG 2004

L5 STRUCTURE UPLOADED

L6 2680 SEARCH L5 SSS FUL

FILE 'CAPLUS' ENTERED AT 12:00:46 ON 18 AUG 2004

FILE 'REGISTRY' ENTERED AT 12:08:19 ON 18 AUG 2004

=> dimethylformamide L12 6859 DIMETHYLFORMAMIDE

=> 18 and 112

'PREP' IS NOT A VALID CROSSOVER QUALIFIER FOR L1
Answer sets created in a different file may be field qualified with a limited set of qualifiers. Enter HELP CROSSOVER at an arrow prompt (=>) for specific information.

=> file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 8.72 206.56 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -3.68

FILE 'CAPLUS' ENTERED AT 12:33:56 ON 18 AUG 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 18 and 112

34275 L12

L13 0 L8 AND L12

=> file reg

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 0.46 207.02 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -3.68

FILE 'REGISTRY' ENTERED AT 12:34:29 ON 18 AUG 2004

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TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> e dimethy	lforma	mide/cn			
E1	1	DIMETHYLFLUOROSILYL (TRICHLORO	OGERMYI.) METHAN	E/CN	
E2	1	DIMETHYLFORMALDEHYDE/CN	JOETHIN JIETHIN	L) CIV	
E3		DIMETHYLFORMAMIDE/CN			
E4	1	DIMETHYLFORMAMIDE 2,4-DINITRO	PHENYT.HYDRAZO	NF/CN	
E5	1	DIMETHYLFORMAMIDE AZINE/CN	Z II DIG IZ O	NE) CIV	
E6	1	DIMETHYLFORMAMIDE COMPLEX WIT	TH MOL TODINE	(1.1)/CN	
E7	1	DIMETHYLFORMAMIDE COMPOUND WI	TH CARRON TET	DACTIODIDE (1.1)/C	
		N	THE CANDON INT.	RACILORIDE (1:1)/C	
E8	1	DIMETHYLFORMAMIDE COMPOUND WI	TH SULFUR TRIC	OXIDE (1.1)/CN	
E9	1	DIMETHYLFORMAMIDE COMPOUND WI	TH WATER (1:1) \CN	
E10	1	DIMETHYLFORMAMIDE COMPOUND WI	TH WATER (2:1) / CN	
E11	1	DIMETHYLFORMAMIDE CYCLIC ETHY	LENE ACETAL/C	N	
E12	1	DIMETHYLFORMAMIDE DEMETHYLASE	E/CN	•	
			,		
=> e3	**				
L14	1 DIM	ETHYLFORMAMIDE/CN			
=> file capl					
COST IN U.S.	DOLLA	RS	SINCE FILE	TOTAL	
ENTRY SESSION					
FULL ESTIMAT	ED COS	Γ	4.85	211.87	
- ·					
DISCOUNT AMO	JNTS (1	FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL	
ENTRY SESSION					
CA SUBSCRIBER PRICE 0.00 -3.68					

FILE 'CAPLUS' ENTERED AT 12:34:56 ON 18 AUG 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 114

L15 26787 L14

=> d his

L5

L7

(FILE 'HOME' ENTERED AT 11:57:09 ON 18 AUG 2004)

FILE 'REGISTRY' ENTERED AT 11:57:45 ON 18 AUG 2004 E TRIPHOSGENE/CN

L1 1 E3

FILE 'CAPLUS' ENTERED AT 11:58:08 ON 18 AUG 2004

L2 854 L1

L3 19404 CHLOROFORMATE

L4 141 L2 AND L3

FILE 'CAPLUS' ENTERED AT 12:00:06 ON 18 AUG 2004

FILE 'REGISTRY' ENTERED AT 12:00:13 ON 18 AUG 2004

STRUCTURE UPLOADED

L6 2680 SEARCH L5 SSS FUL

FILE 'CAPLUS' ENTERED AT 12:00:46 ON 18 AUG 2004

1571 L6/PREP

L8 38 L2 AND L7

L9 379582 AMINE

L10 5 L9 AND L8

L11 0 99999FILE REG

FILE 'REGISTRY' ENTERED AT 12:08:19 ON 18 AUG 2004

L12 6859 DIMETHYLFORMAMIDE

FILE 'CAPLUS' ENTERED AT 12:33:56 ON 18 AUG 2004

L13 0 L8 AND L12

FILE 'REGISTRY' ENTERED AT 12:34:29 ON 18 AUG 2004

E DIMETHYLFORMAMIDE/CN

L14 1 E3

FILE 'CAPLUS' ENTERED AT 12:34:56 ON 18 AUG 2004

L15 26787 L14

=> 114 and 18

26787 L14

L16 0 L14 AND L8

=> save temp 18 chlfmtprep/cn

CHLFMTPREP/CN IS NOT A VALID SAVED NAME

Enter the name you wish to use for the saved query, answer set, or L-number list. The name must:

- 1. Begin with a letter.
- 2. Have 1-12 characters,
- 3. Contain only letters (A-Z) and numbers (0-9),
- 4. End with /Q for a query (search profile,

structure, or screen set), /A for an answer set, or /L for an L-number list.

- 5. Not already be in use as a saved name,
- 6. Not be END, SAV, SAVE, SAVED
- 7. Not have the form of an L-number (Lnnn). ENTER NAME OR (END):end

=> save temp 18 chlfmtprep/a
ANSWER SET L8 HAS BEEN SAVED AS 'CHLFMTPREP/A'

=> save temp 17 chlrfmtes/a
ANSWER SET L7 HAS BEEN SAVED AS 'CHLRFMTES/A'

=> save temp all chlfmtsrch/l L# LIST L1-L16 HAS BEEN SAVED AS 'CHLFMTSRCH/L'

=> logoff hold

COST IN U.S. DOLLARS SINCE FILE TOTAL \ ENTRY SESSION FULL ESTIMATED COST 2.30 214.17 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -3.68

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 12:38:10 ON 18 AUG 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 12:53:03 ON 18 AUG 2004 FILE 'CAPLUS' ENTERED AT 12:53:03 ON 18 AUG 2004 COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	2.30	214.17
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-3.68

=> d his '

(FILE 'HOME' ENTERED AT 11:57:09 ON 18 AUG 2004)

FILE 'REGISTRY' ENTERED AT 11:57:45 ON 18 AUG 2004 E TRIPHOSGENE/CN

L1 1 E3

FILE 'CAPLUS' ENTERED AT 11:58:08 ON 18 AUG 2004

L2 854 L1

L3 19404 CHLOROFORMATE

L4 141 L2 AND L3

FILE 'CAPLUS' ENTERED AT 12:00:06 ON 18 AUG 2004

FILE 'REGISTRY' ENTERED AT 12:00:13 ON 18 AUG 2004

L5 STRUCTURE UPLOADED

L6 2680 SEARCH L5 SSS FUL

FILE 'CAPLUS' ENTERED AT 12:00:46 ON 18 AUG 2004

L7 1571 L6/PREP

L8 38 L2 AND L7

L9 379582 AMINE

L10 5 L9 AND L8

L11 0 99999FILE REG

FILE 'REGISTRY' ENTERED AT 12:08:19 ON 18 AUG 2004

L12 6859 DIMETHYLFORMAMIDE

FILE 'CAPLUS' ENTERED AT 12:33:56 ON 18 AUG 2004

L13 0 L8 AND L12

FILE 'REGISTRY' ENTERED AT 12:34:29 ON 18 AUG 2004

E DIMETHYLFORMAMIDE/CN

L14 1 E3

FILE 'CAPLUS' ENTERED AT 12:34:56 ON 18 AUG 2004

L15 26787 L14

FULL ESTIMATED COST

L16 0 L14 AND L8

SAVE TEMP L8 CHLFMTPREP/A SAVE TEMP L7 CHLRFMTES/A SAVE TEMP ALL CHLFMTSRCH/L

=> file req

COST IN U.S. DOLLARS SINCE FILE TOTAL

ENTRY SESSION 2.75 214.62

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION

CA SUBSCRIBER PRICE ENTRY SESSION 0.00 -3.68

FILE 'REGISTRY' ENTERED AT 12:53:21 ON 18 AUG 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 American Chemical Society (ACS)

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TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

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Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> e triethylamine/cn TRIETHYLALUMINUM-TRIETHYLAMINE COMPLEX/CN 1 E2 1 TRIETHYLALUMINUM-WATER COPOLYMER/CN E3 --> TRIETHYLAMINE/CN TRIETHYLAMINE (2,6-DICHLOROBENZYL)DITHIOCARBAMATE/CN E4 E5 1 TRIETHYLAMINE 1:1 COMPLEX WITH DICHLOROETHYLGERMANE/CN E6 1 TRIETHYLAMINE 1:1 COMPLEX WITH DICHLOROMETHYLGERMANE/CN E7 1 TRIETHYLAMINE 1:1 COMPLEX WITH DICHLOROPHENYLGERMANE/CN TRIETHYLAMINE 2,4,6-TRINITROPHENOLATE/CN F.8 1 TRIETHYLAMINE 2,4-DINITRONAPHTHOLATE/CN E9 1 TRIETHYLAMINE 2,4-DINITROPHENOLATE/CN E10 1 TRIETHYLAMINE 2,6-DICHLOROBENZOATE/CN E111 TRIETHYLAMINE 3,5-DINITROBENZOATE/CN E12 1 => e31 TRIETHYLAMINE/CN L17 => file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION 219.47 4.85

FULL ESTIMATED COST 4.85 219.47

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -3.68

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FILE COVERS 1907 - 18 Aug 2004 VOL 141 ISS 8 FILE LAST UPDATED: 17 Aug 2004 (20040817/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 117 L18 20615 L17

=> d 1129 ti fbib abs L129 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> d 19 ti fbib abs

- L9 ANSWER 1 OF 379582 CAPLUS COPYRIGHT 2004 ACS on STN
- TI Amine organoborane complex initiated polymerizable compositions containing siloxane polymerizable components
- AN 2004:669773 CAPLUS
- TI Amine organoborane complex initiated polymerizable compositions containing siloxane polymerizable components
- IN Sonnenschein, Mark F.; Webb, Steven P.; Wendt, Benjamin L.; Harrington, Daniel R.
- PA Dow Global Technologies Inc., USA
- SO U.S., 20 pp. CODEN: USXXAM
- DT Patent
- LA English

FAN.CNT 1

the

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ΡI	บร 6777512	B1	20040817	US 2003-377440	20030228
				US 2003-377440	20030228

AB In one embodiment the invention is a polymerizable composition comprising a) an organoborane amine complex; b) one or more of monomers,

oligomers or polymers having olefinic unsath, which is capable of polymerization $% \left(1\right) =\left(1\right) +\left(1\right) +\left$

by free radical polymerization; c) one or more compds., oligomers or prepolymers

having a siloxane backbone and reactive moieties capable of polymerization; and d) a catalyst for the polymerization of the one or more compds., oligomers or prepolymers having a siloxane backbone and reactive moieties capable of polymerization. This composition may further comprise a compound which causes

organoborane amine complex to disassoc. In a preferred embodiment, the two part composition further comprises a compound which is reactive with both the b) one or more of monomers, oligomers or polymers having olefinic unsatn. which is capable of polymerization by free radical polymerization; and the c) one or more compds., oligomers or prepolymers having a

siloxane backbone and reactive moieties capable of polymerization This composition

can be polymerized by contacting the two parts of the composition In another embodiment the invention is an organoborane amine complex comprising an alkyl borane having ligands which are alkyl, cycloalkyl or both and an amino siloxane.

=> logoff hold		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	3.60	223.07
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.74	-4.42

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 12:55:06 ON 18 AUG 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 13:17:43 ON 18 AUG 2004 FILE 'CAPLUS' ENTERED AT 13:17:43 ON 18 AUG 2004 COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.60	223.07
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.74	-4.42
=> logoff hold COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY 3.60	SESSION 223.07
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.74	-4.42

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 13:17:50 ON 18 AUG 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1623PAZ

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

Welcome to STN International NEWS 1 Web Page URLs for STN Seminar Schedule - N. America NEWS 2 "Ask CAS" for self-help around the clock NEWS 3 May 12 EXTEND option available in structure searching NEWS Polymer links for the POLYLINK command completed in REGISTRY 4 May 12 NEWS 5 May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus NEWS 6 May 27 CAplus super roles and document types searchable in REGISTRY NEWS 7 Jun 28 Additional enzyme-catalyzed reactions added to CASREACT NEWS Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R) BEILSTEIN enhanced with new display and select options, NEWS 9 Jul 12 resulting in a closer connection to BABS NEWS 10 Jul 30 BEILSTEIN on STN workshop to be held August 24 in conjunction with the 228th ACS National Meeting NEWS 11 AUG 02 IFIPAT/IFIUDB/IFICDB reloaded with new search and display fields AUG 02 CAplus and CA patent records enhanced with European and Japan NEWS 12 Patent Office Classifications NEWS 13 STN User Update to be held August 22 in conjunction with the AUG 02 228th ACS National Meeting NEWS 14 AUG 02 The Analysis Edition of STN Express with Discover!

(Version 7.01 for Windows) now available

NEWS 15 AUG 04 Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004

NEWS 16 AUG 27 BIOCOMMERCE: Changes and enhancements to content coverage

NEWS 17 AUG 27 BIOTECHABS/BIOTECHDS: Two new display fields added for legal status data from INPADOC

NEWS EXPRESS

JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004

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FILE 'HOME' ENTERED AT 07:39:23 ON 31 AUG 2004

=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 07:39:48 ON 31 AUG 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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STRUCTURE FILE UPDATES: 29 AUG 2004 HIGHEST RN 735258-95-4 DICTIONARY FILE UPDATES: 29 AUG 2004 HIGHEST RN 735258-95-4

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Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=> e triphosgene/cn

E1 1 TRIPHOS (FEED SUPPLEMENT)/CN

E2 1 TRIPHOS-I/CN

E3 1 --> TRIPHOSGENE/CN

E4 1 TRIPHOSGENE-PHENYLPHOSPHONIC DICHLORIDE-1,1-BIS(4-HYDROXYPHE NYL)CYCLOHEXANE COPOLYMER/CN

```
1
                   TRIPHOSPHABORETANE/CN
E.6
             1
                   TRIPHOSPHADEN/CN
F.7
             1
                   TRIPHOSPHAGERMETANE/CN
F.8
             1
                   TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIETH
E9
             1
                   TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIPHE
                   NYL-/CN
E10
             1
                   TRIPHOSPHAGERMETANE, 1,2,3-TRIS(1,1-DIMETHYLETHYL)-4,4-DIPHE
                   NYL-, (1A, 2B, 3A) - /CN
E11
                   TRIPHOSPHAGERMETANE, 4,4-DICHLORO-1,2,3-TRIS(1,1-DIMETHYLETH
             1
                   YL)-/CN
E12
             1
                   TRIPHOSPHAMIDE/CN
=> e3
             1 TRIPHOSGENE/CN
=> d 11
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN
     32315-10-9 REGISTRY
CN
     Methanol, trichloro-, carbonate (2:1) (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Carbonic acid, bis(trichloromethyl) ester (6CI, 8CI)
OTHER NAMES:
     Bis(trichloromethyl) carbonate
     Triphosgene
CN
FS
     3D CONCORD
     C3 C16 O3
MF
     COM
CI
LC.
     STN Files:
                  BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS, CASREACT,
       CEN, CHEMCATS, CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, GMELIN*,
       HODOC*, MEDLINE, MSDS-OHS, PROMT, SYNTHLINE, TOXCENTER, USPAT2,
       USPATFULL
         (*File contains numerically searchable property data)
     Other Sources: EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
      CAplus document type: Conference; Journal; Patent
RL.P
       Roles from patents: BIOL (Biological study); PREP (Preparation); PROC
       (Process); RACT (Reactant or reagent); USES (Uses)
       Roles for non-specific derivatives from patents: PREP (Preparation);
RLD.P
       PROC (Process); RACT (Reactant or reagent); USES (Uses)
       Roles from non-patents: ANST (Analytical study); FORM (Formation,
       nonpreparative); MSC (Miscellaneous); OCCU (Occurrence); PREP
       (Preparation); PROC (Process); PRP (Properties); RACT (Reactant or
       reagent); USES (Uses)
RLD.NP Roles for non-specific derivatives from non-patents: PREP
       (Preparation); PRP (Properties); RACT (Reactant or reagent)
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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

856 REFERENCES IN FILE CA (1907 TO DATE)

10 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

858 REFERENCES IN FILE CAPLUS (1907 TO DATE)

3 REFERENCES IN FILE CAOLD (PRIOR TO 1967)